Amendments To The Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

In the Claims:

- 1. (Cancelled).
- 2. (Currently Amended) A compound or salt thereof selected from the group consisting of

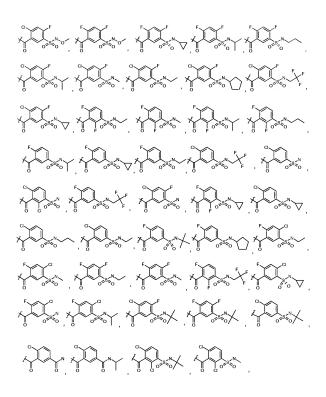
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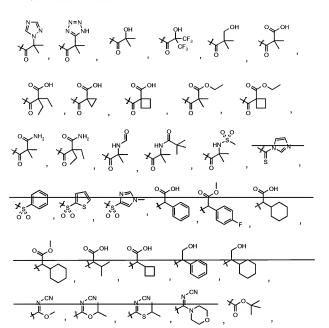
- 3. (Currently amended) A compound selected from the group consisting of tert-butyl 3-(3,4-dichlorophenyl)-3-[3-(4-oxo-1-phenyl-1,3,8-triazaspiro[4.5]dec-8-yl)propyl]pyrrolidine-1-carboxylate;
- 8-{3-[3-(3,4-dichlorophenyl)-1-(2-furoyl)pyrrolidin-3-yl]propyl}-1-phenyl-1,3,8-triazaspiro[4.5]decan-4-one;
- 8-{3-{3-(3,4-dichlorophenyl)-1-(isoxazol-5-ylcarbonyl)pyrrolidin-3-yl]propyl}-1-phenyl-1,3,8-triazaspiro[4.5]decan-4-one;

- 8-{3-{3-(3,4-dichlorophenyl)-1-(1H-pyrrol-2-ylcarbonyl)pyrrolidin-3-yl]propyl}-1-phenyl-1.3.8-triazaspirol4.5ldecan-4-one:
- $8-\{3-[3-(3,4-dichlorophenyl)-1-pentanoylpyrrolidin-3-yl]propyl\}-1-[3-(3,4-dichlorophenyl)-1-[3$
- (trifluoromethyl)phenyl]-1,3,8-triazaspiro[4.5]decan-4-one;
- 8-{3-[3-(3,4-dichlorophenyl)-1-(2-furoyl)pyrrolidin-3-yl]propyl}-1-[3-
- (trifluoromethyl)phenyl]-1,3,8-triazaspiro[4.5]decan-4-one;
- 8-{3-[1-(cyclobutylcarbonyl)-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]propyl}-1-[3-
- (trifluoromethyl)phenyl]-1,3,8-triazaspiro[4.5]decan-4-one;
- 8-{3-[3-(3,4-dichlorophenyl)-1-pentanoylpyrrolidin-3-yl]propyl}-1-phenyl-1,3,8-triazaspirol4.5]decan-4-one:
- 8-{3-[1-(cyclopentylcarbonyl)-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]propyl}-1-phenyl-
- 1,3,8-triazaspiro[4.5]decan-4-one;
- 8-{3-[1-(cyclobutylcarbonyl)-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]propyl}-1-phenyl-1,3,8-triazaspiro[4.5]decan-4-one;
- 8-{3-[1-(cyclobutylcarbonyl)-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]propyl}-1-(3-methylphenyl)-1,3,8-triazaspiro[4.5]decan-4-one;
- 3-acetyl-8-{3-[1-acetyl-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]propyl}-1-(3-methylphenyl)-1.3.8-triazaspirol4.5ldecan-4-one:
- 8-{3-[1-(1,3-benzoxazol-2-yl)-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]propyl}-1-phenyl-
- 1,3,8-triazaspiro[4.5]decan-4-one;
- 8-(2-[[3-(3,4-dichlorophenyl)-1-(2-furoyl)pyrrolidin-3-yl]oxy}ethyl)-1-phenyl-1,3,8-triazaspirol4.5]decan-4-one:
- 8-(2-[[1-(cyclopentylcarbonyl)-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]oxy}ethyl)-1-phenyl-1.3.8-triazaspiro[4.5]decan-4-one:
- 8-(2-{[1-acetyl-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]oxy}ethyl)-1-phenyl-1,3,8-triazaspirol4.5ldecan-4-one:
- 8-(2-[[3 (3,4-dichlorophenyl)-1 (phenylsulfonyl)pyrrolidin-3-yl]oxy}ethyl) 1-phenyl-1,3,8-triazaspiro[4-5]decan-4-one;
- 8-(2-[[3-(3,4-dichlorophenyl)-1-(2-furoyl)pyrrolidin-3-yl]oxy}ethyl)-1-(3-methoxyphenyl)-1,3,8-triazaspiro[4.5]decan-4-one;

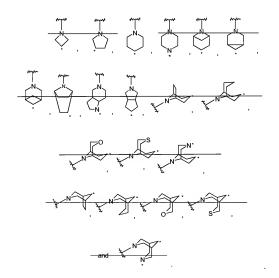
- 8-(2-{[1-(cyclopentylcarbonyl)-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]oxy}ethyl)-1-(3-methoxyphenyl)-1.3.8-triazaspirol4.5ldecan-4-one; and
- 8-(2-[[1-acetyl-3-(3,4-dichlorophenyl)pyrrolidin-3-yl]oxy}ethyl)-1-(3-methoxyphenyl)-1.3.8-triazaspiro[4.5]decan-4-one:-and
- 8-(2-[[3-(3,4-dichlorophenyl)-1-(phenylsulfonyl)pyrrolidin-3-yl]oxy]ethyl)-1-(3-methoxyphenyl)-1,3,8-triazaspiro[4-5]decan-4-one.
- 4. (Currently Amended) The compound of claim 4 <u>48</u> wherein the B ring is pyrrolidine.
- 5. (Original) The compound of claim 4 wherein R⁹ is H.
- (Cancelled).
- (Currently Amended) The compound of claim 4 48 wherein R¹ is anyl or a substituted anyl.
- 8. (Original) The compound of claim 7 wherein R¹ is phenyl mono- or di- substituted with halogen.
- 9. (Original) The compound of claim 8 wherein R¹ is phenyl di-substituted with Cl.
- 10. (Currently Amended) The compound of claim $4 \underline{48}$ wherein $-(Y)_m R^3$ is selected from the group consisting of

11. (Currently Amended) The compound of claim $4\,\underline{48}$ wherein $-(Y)_m$ - R^3 is selected from the group consisting of

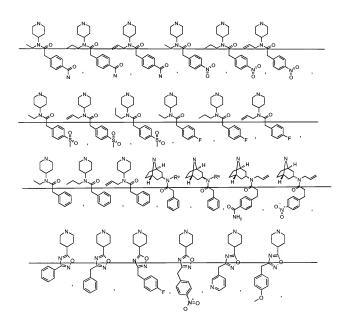


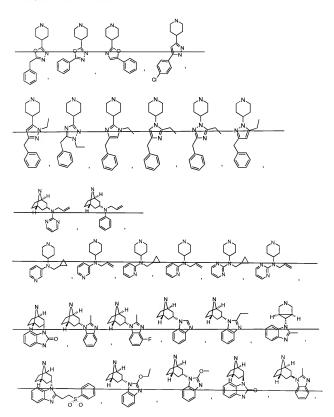


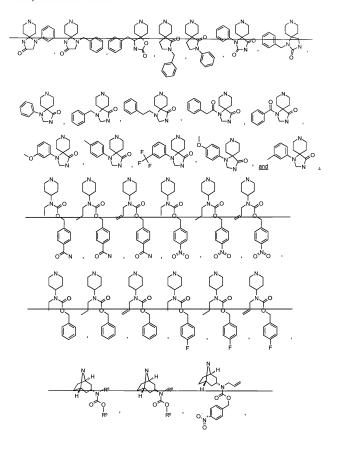
- 12. (Currently Amended) The compound of claim 4.48 wherein m is 1, Y is -C(O), and R^3 is either aryl or heteroaryl, wherein said aryl or heteroaryl is optionally substituted, with an optionally substituted alkyl, or an optionally substituted cycloalkyl.
- 13. (Cancelled).
- 14. (Cancelled).
- 15. (Currently Amended) The compound of claim 4.48 wherein m is 1, Y is -C(O)O, and R^3 is optionally substituted alkyl or optionally substituted aryl.
- 16.-24. (Cancelled).
- (Currently Amended) The compound of claim 4 48 wherein the A ring, with an
 asterisk indicating the point of substitution, is selected from the group consisting of

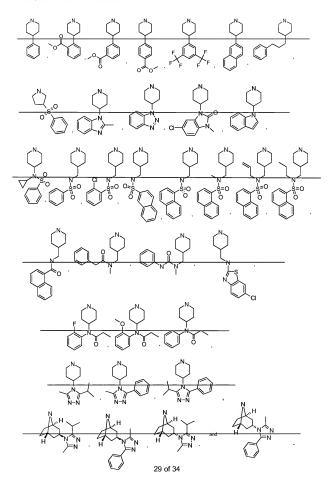


- 26.- 28. (Cancelled).
- 29. (Currently amended) The compound of claim 27 48 wherein the A ring ring in combination with R² is selected from the group consisting of[[.]]









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30.- 42. (Cancelled).

- (Currently Amended) A pharmaceutical composition comprising a
 pharmaceutically effective amount of a compound according to elaims 1-30 claim
 48 together with a pharmaceutically acceptable carrier.
- (Currently Amended) The pharmaceutical composition according to claim 42 43 in the form of a tablet or capsule.
- (Currently Amended) The pharmaceutical composition according to claim 42 43in the form of a liquid.
- 46. 47. (Cancelled).
- 48. (New) A compound of formula (I)

and pharmaceutically acceptable derivatives thereof, wherein

X is (CH₂)₃;

Ring A is a saturated, 6-membered monocyclic ring having one ring nitrogen; Ring B is a saturated 4 or 5 membered ring containing the depicted ring nitrogen; R^1 is a phenyl or a phenyl substituted with mono- or di- halogen;

Y is -C(O)- or -C(O)-O, and m is 1;

 $R^{3} \text{ is H, -N(R^{0})}_{2}, -N(R^{0})C(O)R^{0}, -CN, \text{ halogen, CF}_{3}, \text{ alkyl optionally substituted by one or more groups selected from R^{7} or -S-aryl optionally substituted by -(CH₂)₁₋₆-N(R^{0})SO₂(R^{0}), \text{ alkenyl optionally substituted by one or more groups selected from R^{7} or -S-aryl optionally substituted by -(CH₂)₁₋₆-N(R^{0})SO₂(R^{0}), \text{ alkynyl optionally substituted by one or more groups selected from R^{7} or -S-aryl optionally substituted by -(CH₂)₁₋₆-N(R^{0})SO₂(R^{0}), cycloalkyl or carbocyclyl optionally substituted by one or more R^{6}, heteroaryl optionally substituted by one or more R^{6}, or heterocyclyl optionally substituted by one or more R^{6};$

each R^6 is independently selected from the group consisting of halogen, -CF3, -OCF3,

 $-\mathsf{OR}^0, -(\mathsf{CH}_2)_{1:6} - \mathsf{OR}^0, -\mathsf{SR}^0, -(\mathsf{CH}_2)_{1:6} - \mathsf{SR}^0, -\mathsf{SCF}_3, -\mathsf{R}^0, \text{ methylenedioxy},\\ \text{ethylenedioxy}, -\mathsf{NO}_2, -\mathsf{CN}, -(\mathsf{CH}_2)_{1:6} - \mathsf{CN}, -\mathsf{N}(\mathsf{R}^0)_2, -(\mathsf{CH}_2)_{1:6} - \mathsf{N}(\mathsf{R}^0)_2, -\mathsf{NR}^0 - \mathsf{C}(\mathsf{Q})_{\mathsf{N}}^0,\\ -\mathsf{NR}^0(\mathsf{CN}), -\mathsf{NR}^0\mathsf{C}(\mathsf{Q})\mathsf{N}(\mathsf{R}^0)_2, -\mathsf{NR}^0\mathsf{C}(\mathsf{Q})\mathsf{N}(\mathsf{R}^0)_2, -\mathsf{NR}^0\mathsf{NR}^0\mathsf{C}(\mathsf{Q})\mathsf{R}^0,\\ -\mathsf{NR}^0\mathsf{NR}^0\mathsf{C}(\mathsf{Q})\mathsf{N}(\mathsf{R}^0)_2, -\mathsf{NR}^0\mathsf{NR}^0\mathsf{C}\mathsf{Q}_2\mathsf{R}^0, -\mathsf{C}(\mathsf{Q})\mathsf{C}(\mathsf{Q})^0,\\ -(\mathsf{CH}_2)_{0:6}\mathsf{CO}_2\mathsf{R}^0, -\mathsf{C}(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0,\\ -(\mathsf{C}(\mathsf{H}_2)_{0:6}\mathsf{C}^0, -\mathsf{C}(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0,\\ -(\mathsf{C}(\mathsf{Q})\mathsf{N}(\mathsf{R}^0)_2, -\mathsf{C}(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0,\\ -(\mathsf{Q})^0, -(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0,\\ -(\mathsf{Q})^0, -(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0,\\ -(\mathsf{Q})^0, -(\mathsf{Q})^0, -\mathsf{C}(\mathsf{Q})^0,\\ -(\mathsf{Q})^0, -(\mathsf{C})^0, -(\mathsf{Q})^0, -(\mathsf{Q})^0,\\ -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0,\\ -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0,\\ -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0,\\ -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0, -(\mathsf{C})^0,\\ -(\mathsf{C})^0, -$

each R^7 is independently selected from the group consisting of halogen, $-CF_3$, $-R^0$, $-OR^0$, $-OCF_3$, $-(CH_2)_{1-6}-OR^0$, $-SR^0$, $-SCF_3$, $-(CH_2)_{1-6}-SR^0$, aryl optionally substituted by R^6 , methylenedioxy, ethylenedioxy, $-NO_2$, -CN, $-(CH_2)_{1-6}-CN$, $-N(R^0)_2$, $-(CH_2)_{1-6}-N(R^0)_2$, $-NR^0C(O)R^0$, $-NR^0(CN)$, $-NR^0C(O)N(R^0)_2$, $-NR^0NR^0C(O)R^0$, $-NR^0NR^0C(O)R^0$, $-NR^0NR^0C(O)R^0$, $-NR^0NR^0CO_2R^0$, $-C(O)C(O)R^0$, $-C(O)CH_2C(O)R^0$, $-(CH_2)_{1-6}-CO_2R^0$, $-C(O)R^0$, $-C(O)N(R^0)N(R^0)_2$, $-C(O)N(R^0)N(R^0)CO_2R^0$, $-C(O)N(R^0)CO_2R^0$, $-C(O)N(R^0)N(R^0)CO_2R^0$,

each R^8 is independently selected from R^7 , =0, =S, =N(R^0), and =N(CN); R^9 is H:

each R⁰ is independently selected from the group consisting of hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl, carbocyclylalkyl, aryl, heteroaryl, aralkyl, heteroaralkyl, heterocyclyl, and heterocyclylalkyl, wherein each member of R⁰ except H is

optionally substituted by one or more R*, OR*, N(R*)₂, =O, =S, halo, CF₃, NO₂, CN, -C(O)R*, -CO₂R*, -C(O)-aryl, -C(O)-heteroaryl, -C(O)-aralkyl, -S(O),-aryl, -S(O)₁-heteroaryl, -NR*SO₂R*, -NR*C(O)R*, -NR*C(O)N(R*)₂, -N(R*)C(S)N(R*)₂, -NR*NR*C(O)R*, -NR*NR*C(O)N(R*)₂, -NR*NR*CO₂R*, -C(O)C(O)R*, -C(O)CH₂C(O)R*, -C(O)N(R*)N(R*)₂, -C(O)N(R*)₂, -C(O)N(R*)₂, -C(O)N(R*)₂, -SO₂N(R*)₂ wherein the two R*s on the same nitrogen are optionally taken together to form a 5-8 membered saturated, partially saturated or aromatic ring having additional 0-4 heteroatoms selected from oxygen, phosphorus, nitrogen or sulfur; each R* is independently H. alkyl, alkenyl, alkynyl, cycloalkyl, aryl, or heteroaryl.